

# Implicit Methods and Globalization Strategies for the Robust Approximation of Solutions to the Reynolds Averaged Navier-Stokes equations

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Knowledge for Tomorrow

# Outline

- Introduction and Motivation – The RANS equations
- Solution algorithms – Multigrid smoothers
- Globalization strategies
- Numerical examples



# Motivation

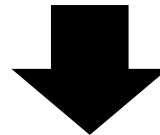
## Goals of flighpath 2050:

1. In 2050 technologies and procedures available allow a 75% reduction in CO<sub>2</sub> emissions per passenger kilometre to support the ATAG (Air Transport Action Group) target and a 90% reduction in NO<sub>x</sub> emissions.
2. The perceived noise emission of flying aircraft is reduced by 65%.

These are relative to the capabilities of typical new aircraft in 2000.

3. Overall, the European air transport system has less than one accident per ten million commercial aircraft flights.

→ The future aircraft is ecologically sensitive, low noise, and safe.



A key element, to design aircrafts ready for the future, is the **accurate and efficient simulation of fluid flow** coupled with other disciplines such as aerodynamics and aeroacoustics.



# Requirements of a CFD code

- Reliable tool in a process chain
- Interaction with other components (e.g. structure, mesh deformation, ...)
- Accuracy, e.g. prediction of force coefficients up to a certain accuracy
- Evaluation and assessment of turbulence models
- ...

## Basic demand:

- Machine accurate solutions (on a given grid, that is a given resolution)
- Mesh converged solutions (in general hard to obtain, in particular in 3D)
- The code needs to run on regular basis without user interaction



# How to prepare a CFD code for the future

- Identify the main building blocks
- Disaggregate the code into its building blocks
- Write routines which check the building blocks with respect to correctness
- Design interfaces such that the building blocks can be easily exchanged



## Modular software design

**Requirement: Identification of the main building blocks of a CFD code, to  
create algorithms prepared for demands in fully automatic process  
chains.**



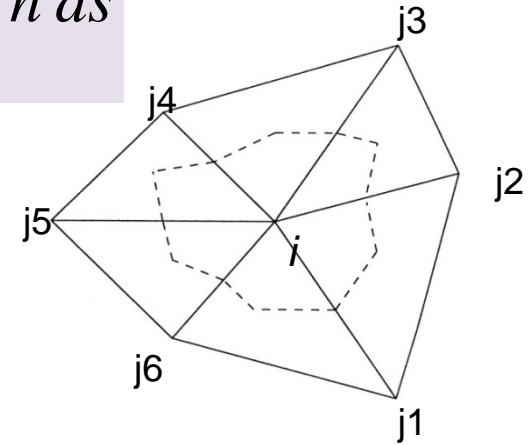
# Compressible Navier-Stokes equations

$W := (\rho, \rho u, \rho E) \rightarrow$  Conservative variables

$$\frac{d}{dt} \int_{\Omega} W dx = - \int_{\partial\Omega} (F_c(W, \text{grad } W) - F_v(W, \text{grad } W)) \cdot n ds$$

$F_c$  : Convective flux

$F_v$  : Viscous flux



## Unstructured Finite volume discretization

→ Nonlinear operator equation:

$$\frac{dW}{dt} = -M^{-1}R(W), \quad M = \text{diag}(\text{vol}(\Omega_i)_{i=1,\dots,N})$$

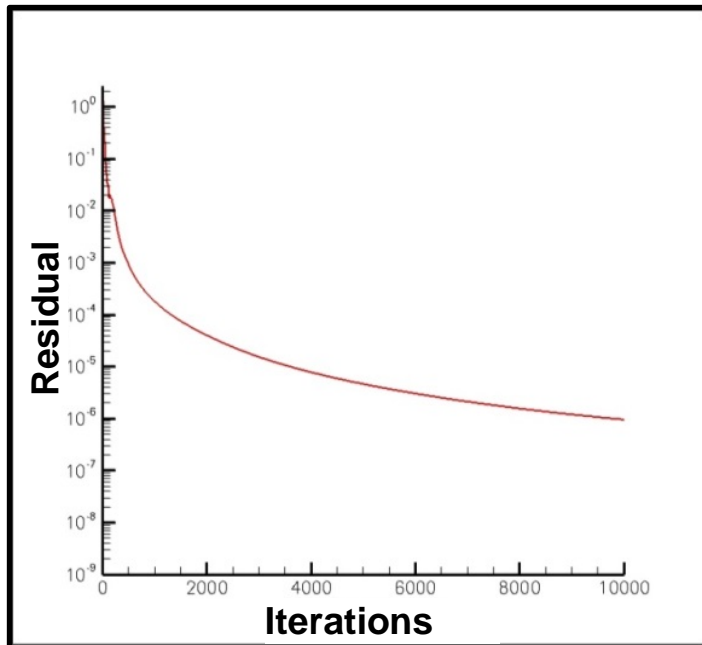
Interested in steady state solution:  $R(W) = 0$



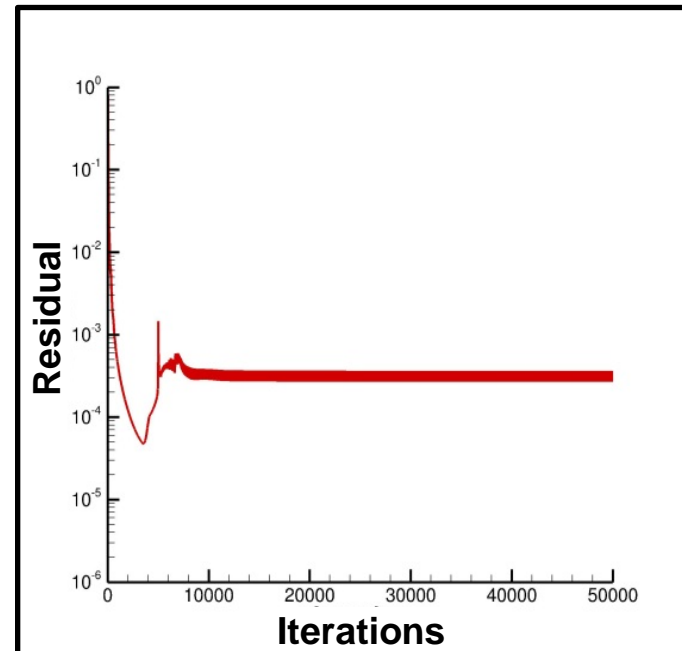


# Necessity for improvement of solution algorithms

## Typical convergence behavior for high Reynolds number viscous flows

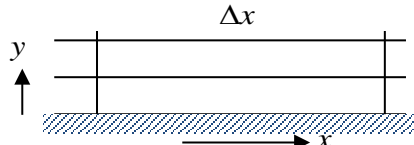


➡ Convergence rate deteriorates significantly after initial phase



➡ No convergence, iteration stagnates

➔ Stiff set of equations



A diagram of a rectangular domain with horizontal dimension  $\Delta x$  and vertical dimension  $\Delta y$ . The x-axis is horizontal and the y-axis is vertical. The domain is divided into a grid of cells. The aspect ratio is given by the equation:

$$\frac{\Delta x}{\Delta y} > 10000$$

- anisotropic cells to represent gradients in the boundary layer
- turbulent flow equations with source terms



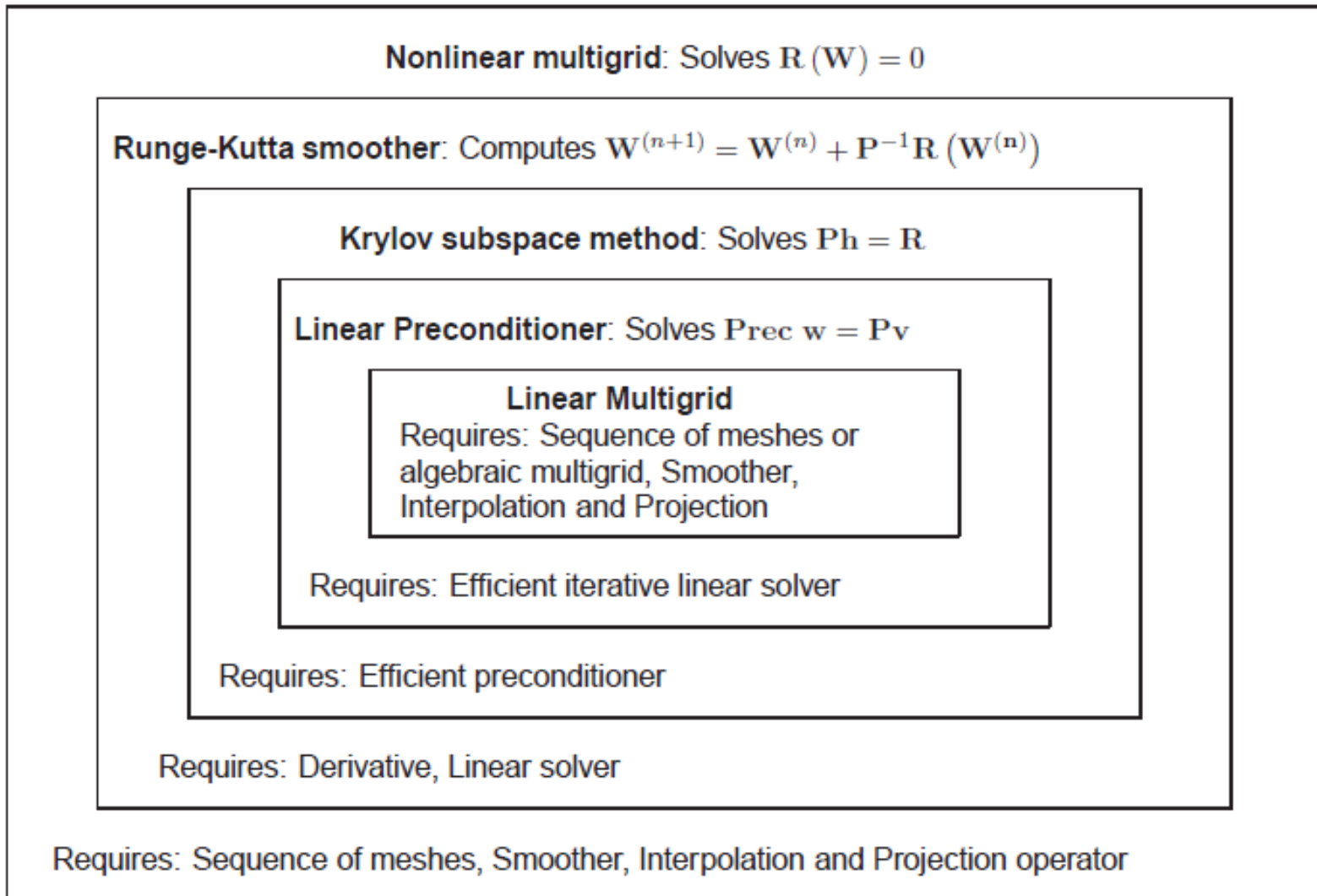
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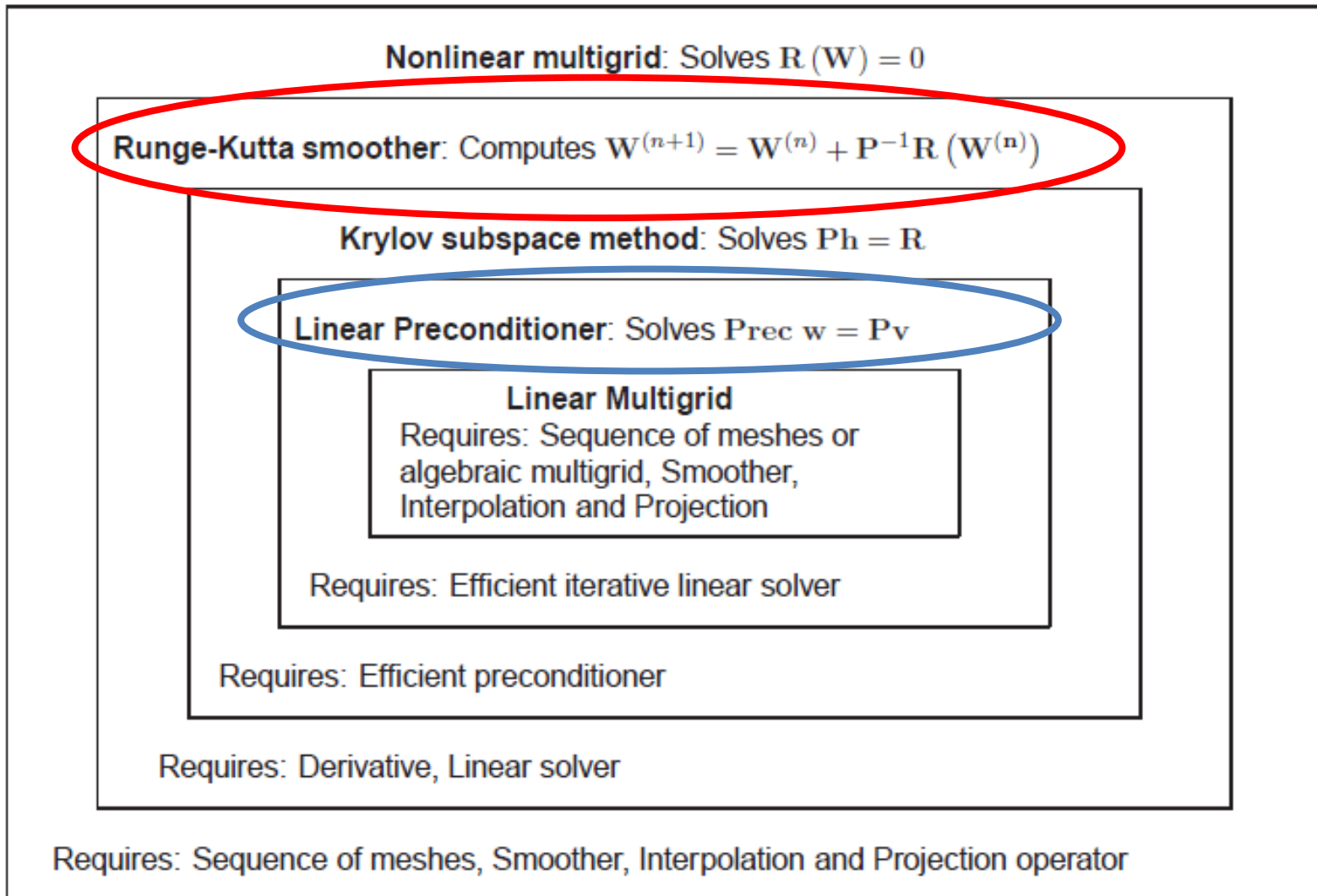




# Structure of solution algorithm



# Structure of solution algorithm



# Multigrid smoother: **Prec.** Runge-Kutta method

$$\frac{dW}{dt} = -M^{-1}R(W), \quad M = \text{diag}(\text{vol}(\Omega_i)_{i=1,\dots,N})$$

Apply **preconditioned** expl. Runge-Kutta method

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} P_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s \\ W_{n+1} &:= W^{(s)} \\ P_j &:= (\Delta t)^{-1} M + \frac{\partial R}{\partial W} \end{aligned}$$

to approximate  $W$  such that  $R(W) \approx 0$



# Multigrid smoother: **Prec.** Runge-Kutta method

$$\frac{dW}{dt} = -M^{-1}R(W), \quad M = \text{diag}(\text{vol}(\Omega_i)_{i=1,\dots,N})$$

Apply **preconditioned** expl. Runge-Kutta method

$$W^{(0)} := W_n$$

$$W^{(j)} := W^{(0)} - \alpha_{j+1,j} \left( \frac{CFL \Delta t_i}{\text{vol}(\Omega_i)} \right) R(W^{(j-1)}), \quad j = 1, \dots, s$$

$$W_{n+1} := W^{(s)}$$

$$P_j := (\Delta t)^{-1} M + \cancel{\frac{\partial R}{\partial W}} = \left( \frac{CFL \Delta t_i}{\text{vol}(\Omega_i)} \right)^{-1} = \left( \frac{\text{vol}(\Omega_i)}{CFL \Delta t_i} \right)^{-1}$$

Explicit Runge  
Kutta scheme with  
local time stepping

to approximate  $W$  such that  $R(W) \approx 0$

→ Requires inversion of a scalar value for each control volume:

$$\left( \frac{\text{vol}(\Omega_i)}{CFL \Delta t_i} \right)^{-1}$$



# Multigrid smoother: **Prec.** Runge-Kutta method

$$\frac{dW}{dt} = -M^{-1}R(W), \quad M = \text{diag}(\text{vol}(\Omega_i)_{i=1,\dots,N})$$

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$$x = P_j^{-1} R(W^{(j-1)}) \Leftrightarrow P_j x = R(W^{(j-1)})$$

to approximate  $W$  such that  $R(W) \approx 0$

→ Task: Need to approximate efficiently solution of  $Px = R(W)$

→ Inversion of scalar value is replaced by solving a large scale linear system.



# The connection to Newton's method

Outer Loop: Multistage Runge-Kutta method  $\rightarrow$  Choose  $s = 1$ , i.e. only one stage

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} P_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s \\ W_{n+1} &:= W^{(s)} \end{aligned}$$

$$P_j := (\Delta t)^{-1} M + \frac{\partial R}{\partial W} = \frac{\partial R}{\partial W}, \quad CFL \rightarrow \infty$$

$$(\Delta t)^{-1} = \left( \frac{\text{vol}(\Omega_i)}{CFL \Delta t_i} \right) \rightarrow 0, \quad CFL \rightarrow \infty$$

$$W_{n+1} := W_n - \left( \frac{\partial R}{\partial W} \right)^{-1} R(W_n)$$

$\rightarrow$  The solution method is some kind of generalization of Newton's method

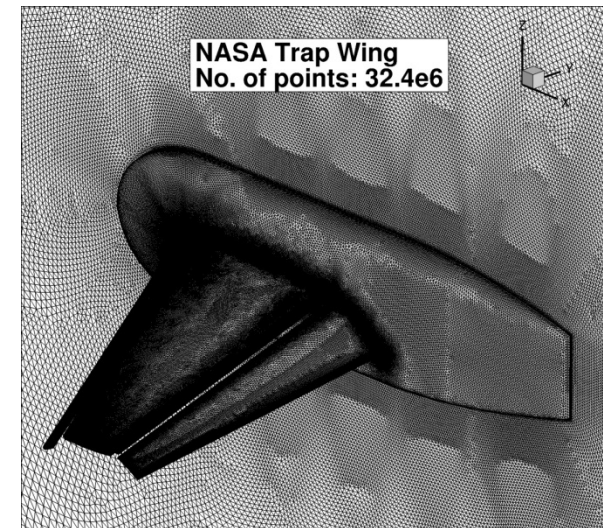




$$P_x = R(W)$$

- Represents a large scale (in general more than  $10^8$  unknowns), ill-conditioned linear system
- It is not of interest to solve these linear systems, it is of interest to get a ***reasonable update*** for the outer nonlinear loop
- Krylov subspace methods are a natural choice for a matrix-free implementation
- A well suited **preconditioner** is required

**Krylov subspace methods are in general only effective in combination with a well suited preconditioner!**





# Code design

- execute multigrid cycle (until convergence)
  - execute preconditioned expl. Runge-Kutta algorithm
    - evaluate residual  $R$
    - evaluate derivative  $dR$ ,  $P = \Delta t + dR$
    - solve linear system  $(P, R)$ 
      - apply preconditioned Krylov subspace method
        - construct a (further) preconditioner for the linear system  $Px = R$
- update flow variables  $W$

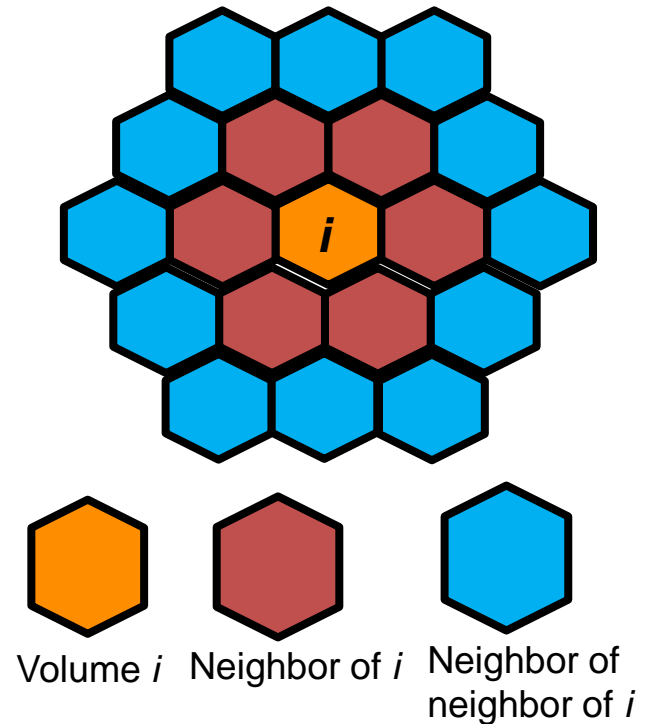


# Construction of Preconditioner (for lin. System)

Idea: Base preconditioner upon next neighbor stencil

$$R_i^{2nd}(W_i, W_{j \in N(i)}, W_{k, k \in N(j)}) \approx R_i^{1st}(W_i, W_{j, j \in N(i)})$$

$$\Rightarrow \frac{\partial R^{2nd}}{\partial W} \approx \frac{\partial R^{1st}}{\partial W} \Rightarrow \text{Prec} = (\Delta t)^{-1} M + \frac{\partial R^{1st}}{\partial W}$$



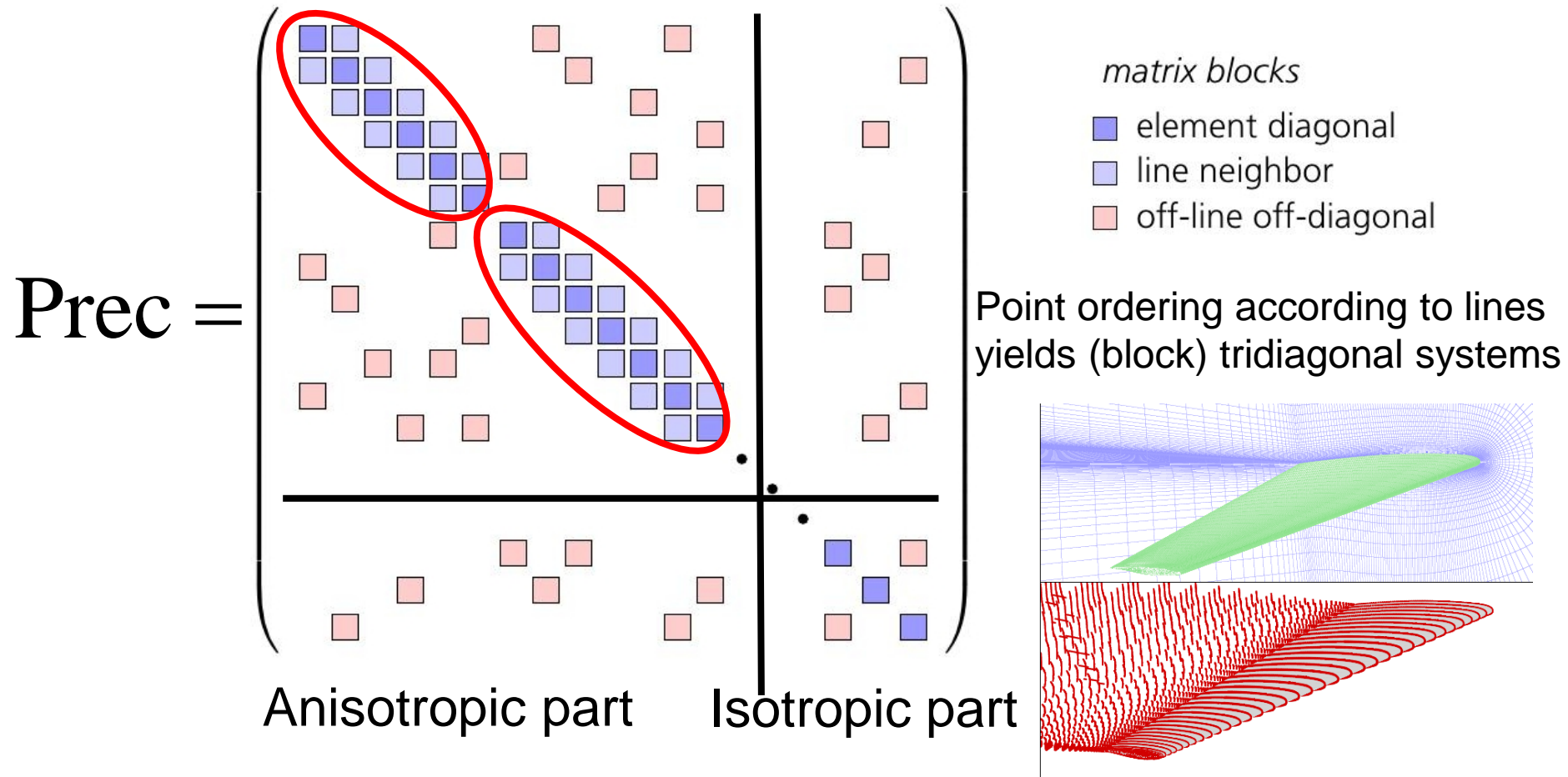
→ Required: Solution method for **Prec  $w = b$**



# Challenge: Find approximate solution of linear system

$$\text{Prec } w = b$$

where Prec is a block sparse matrix of dimension number of mesh points



# Iterative solution methods for $\text{Prec } w = z$

**Mathematical textbook** methods for solution of linear systems, e.g

- (Block-) Jacobi method
- (Block-) Gauss-Seidel method
- Symmetric (Block-) Gauss-Seidel method

Methods have been extended:

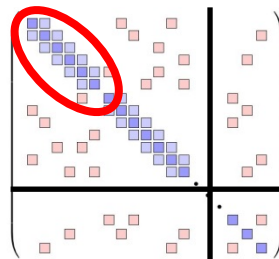
→ Exploit ***directions of strongest coupling*** in iterative solution process

**(Symmetric) Line (Jacobi) Gauss-Seidel method:**

$$x_{L_i}^{(m+1)} = \text{tridiag}(D_{L_i})^{-1} \left( b_{L_i} - \sum_{j \in L_1, \dots, L_{i-1}, j \notin L_i} \text{Prec}_{L_i j} x_j^{(m+1)} - \sum_{j \notin L_1, \dots, L_{i-1}, j \notin L_i} \text{Prec}_{L_i j} x_j^{(m)} \right)$$



**Algebraic representation  
and implementation of  
geometric data (Lines)**



# Code design

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    - evaluate residual  $R$
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    - solve linear system  $(P, R)$ 
      - apply preconditioned Krylov subspace method
        - construct a (further) preconditioner for the linear system  $Px = R$
        - solve  $P_{\text{prec}} w = b$  to precondition  $Px = R$  (by Line symmetric Gauss-Seidel method)
  - update flow variables  $W$



# A historical view on solution methods in CFD

## Two competitive views

Multigrid + Low cost smoother  $\leftrightarrow$  Newton's method (expensive smoother)

### Low cost smoothers:

1. Expl. Runge-Kutta + local time stepping (Jameson)
2. Point implicit Runge-Kutta (Pierce, Giles, Moinier)
3. Line implicit Runge-Kutta (Mavriplis)
4. 1.st order approximate Jacobian (Swanson, Rossow, Yoon + Jameson (LU-SGS))

### Preconditioned explicit Runge-Kutta smoother

All well known specific smoothers developed throughout the CFD literature are specifications of the general method shown here

The suggested methods just differ with respect to the ***approximation of the exact Jacobian*** and the ***iterative solver***





# Derivation of low cost smoothers

$$W^{(0)} := W_n$$

$$W^{(j)} := W^{(0)} - \alpha_{j+1,j} P_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s$$

$$W_{n+1} := W^{(s)}$$

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR$ ,  $P = \Delta t + dR$
  - **solve linear system  $(P, R)$** 
    - **apply preconditioned Krylov subspace method**
      - construct a (further) preconditioner for the linear system  $Px = R$
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$$W_{n+1} := W^{(s)}$$

Simplifications:

1. Number of Krylov steps = 0

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual R
  - evaluate derivative dR,  $P = \Delta t + dR$
  - ~~solve linear system (P, R)~~
    - ~~apply preconditioned Krylov subspace method~~
      - construct a (further) preconditioner for the linear system  $Px = R$
      - solve  $P_{\text{prec}} w = b$  to precondition  $Px = R$
  - update flow variables W



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Simplifications:

1. Number of Krylov steps = 0

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $Prec = \Delta t + dR^{1st}$ 
    - ~~• construct a (further) preconditioner for the linear system  $Px = R$~~
    - ~~• solve  $Prec w = b$  to precondition  $Px = R$~~
  - update flow variables  $W$



# Derivation of low cost smoothers

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} \text{Prec}_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s \\ W_{n+1} &:= W^{(s)} \end{aligned}$$

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## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st}$ 
    - solve  $\text{Prec } w = b$
  - update flow variables  $W$

Preconditioning based on 1.st order approximate Jacobian  
(Swanson, Rossow, Yoon + Jameson (LU-SGS))



# Derivation of low cost smoothers

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} \text{Prec}_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s \\ W_{n+1} &:= W^{(s)} \end{aligned}$$

Simplifications:

1. Number of Krylov steps = 0
2. Simplify  $dR^{1st}$  entries by spectral radius

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st} \approx \Delta t + \varrho(dR^{1st})$ 
    - solve  $\text{Prec } w = b$
  - update flow variables  $W$

Solve with 1 symmetric Gauss-Seidel sweep

Yoon + Jameson (LU-SGS)



# Derivation of low cost smoothers

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} \text{Prec}_j^{-1} R(W^{(j-1)}), \quad j = 1, \dots, s \\ W_{n+1} &:= W^{(s)} \end{aligned}$$

Simplifications:

1. Number of Krylov steps = 0
2. Simplify  $dR^{1st}$  entries by spectral radius
3. Iterative solver: Line Jacobi truncated after one step

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual R
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st}$ 
    - solve  $\text{Prec} w = b$
  - update flow variables W

$$x^{(0)} = 0$$

$$x_{L_i}^{(m+1)} = \text{tridiag}(D_{L_i})^{-1} \left( b_{L_i} - \sum_{j \notin L_1, \dots, L_{i-1}, j \notin L_i} \text{Prec}_{L_i j} x_j^{(m)} \right)$$



# Derivation of low cost smoothers

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- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st}$ 
    - solve  $\text{Prec } w = b$
  - update flow variables  $W$

$$x^{(0)} = 0$$

$$x_{L_i}^{(1)} = \text{tridiag}(D_{L_i})^{-1} b_{L_i}$$



# Derivation of low cost smoothers

$$W^{(0)} := W_n$$

$$W_{L_k}^{(j)} := W_{L_k}^{(0)} - \alpha_{j+1,j} \text{tridiag}(D_{L_k})^{-1} R(W^{(j-1)})$$

$$W_{n+1} := W^{(s)}$$

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st}$ 
    - solve  $w = \text{tridiag}(D)^{-1}b$
  - update flow variables  $W$

**Line implicit Runge-Kutta (Mavriplis)**

Simplifications:

1. Number of Krylov steps = 0
2. Simplify  $dR^{1st}$  entries by spectral radius
3. Iterative solver: Line Jacobi truncated after one step

$$x^{(0)} = 0$$

$$x_{L_i}^{(1)} = \text{tridiag}(D_{L_i})^{-1} b_{L_i}$$





# Derivation of low cost smoothers

$$W^{(0)} := W_n$$

$$W_k^{(j)} := W_k^{(0)} - \alpha_{j+1,k} (D_k)^{-1} R(W^{(j-1)})$$

$$W_{n+1} := W^{(s)}$$

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $Prec = \Delta t + dR^{1st}$ 
    - solve  $w = (D)^{-1}b$
  - update flow variables  $W$

## Point implicit Runge-Kutta (Giles, Moinier)

Simplifications:

1. Number of Krylov steps = 0
2. Simplify  $dR^{1st}$  entries by spectral radius
3. Iterative solver: Line Jacobi truncated after one step
4. Neglect lines, i.e. perform only Jacobi iteration

$$x^{(0)} = 0$$

$$x_i^{(1)} = (D_i)^{-1} b_i$$



# Derivation of low cost smoothers

$$W^{(0)} := W_n$$

$$W_k^{(j)} := W_k^{(0)} - \alpha_{j+1,j} \frac{\Delta t_k}{\text{vol}(\Omega_k)} R(W^{(j-1)})$$

$$W_{n+1} := W^{(s)}$$

## Smoothing step

- execute preconditioned expl. Runge-Kutta algorithm
  - evaluate residual  $R$
  - evaluate derivative  $dR^{1st}$ ,  $\text{Prec} = \Delta t + dR^{1st}$ 
    - solve  $w = (\Delta t / \text{vol})^{-1} b$
  - update flow variables  $W$

**Explicit Runge-Kutta (Jameson)**

**→ Full hierarchy of solution methods**

Simplifications:

1. Number of Krylov steps = 0
2. Simplify  $dR^{1st}$  entries by spectral radius
3. Iterative solver: Line Jacobi truncated after one step
4. Neglect lines, i.e. perform only Jacobi iteration
5. Approximate diagonal terms of Jacobian by spectral radius

$$x_i^{(1)} = (\rho(D_i))^{-1} b_i = \frac{\Delta t_i}{\text{vol}(\Omega)_i} b_i$$



# Main building blocks of a CFD code

- Data structure for (block) sparse matrices  $\rightarrow \frac{\partial R}{\partial W}$
- Data structure for (block) vectors  $\rightarrow R(W)$
- Algorithms acting on these data structures

$\rightarrow$  Link to a suited, efficient **LINEAR ALGEBRA PACKAGE**



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# What is a globalization strategy?

Consider Newton's method  $W^{n+1} = W^n - \left[ \frac{\partial R}{\partial W} \right]^{-1} R(W^n)$

A globalization strategy is the try to construct an algorithm which

1. preserves the nice properties of Newton's method
2. circumvents its shortcomings

## Why do we need it?

Newton's method converges only

1. under certain smoothness assumptions
2. if the initial guess is in a neighborhood of the root
3. ...



**Parameter settings allow for several possible smoothing techniques:**

- Number of Runge-Kutta stages
- Number of Gauss-Seidel sweeps
- Number of Krylov subspace steps
- Approximation of Jacobian
- .....

Newton's method



Explicit Runge-Kutta

## How to choose a robust and efficient method?

→ Development of an analysis tool to give some guideline



# Evaluation of smoother:

## Consider linearized problem

### Nonlinear Problem:

$$\frac{dW}{dt} = -M^{-1}R(W)$$

### Linearized Problem

$$\frac{dW}{dt} \approx -M^{-1} \left( \underbrace{R(W^*)}_{=0; \text{Steady state}} + \frac{\partial R}{\partial W} [W^*] \Delta W \right)$$

$$\begin{aligned} W^{(0)} &:= W_n \\ W^{(j)} &:= W^{(0)} - \alpha_{j+1,j} P_j^{-1, \text{app}} R(W^{(j-1)}) \\ W_{n+1} &:= W^{(s)} \end{aligned}$$

$$W^{(0)} := W_n$$

$$W^{(j)} := W^{(0)} - \alpha_{j+1,j} P_j^{-1, \text{app}} \frac{\partial R}{\partial W} W^{(j-1)}$$

$$W_{n+1} := W^{(s)} \Leftrightarrow W^{(n+1)} = q_s \left( P_j^{-1, \text{app}} \frac{\partial R}{\partial W} \right) W^{(j-1)}$$

$$q_s(z) = 1 + \sum_{j=1}^s \beta_j z^j$$

$$\text{Convergence} \Leftrightarrow \rho \left( q_s \left( P_j^{-1, \text{app}} \frac{\partial R}{\partial W} \right) \right) < 1$$





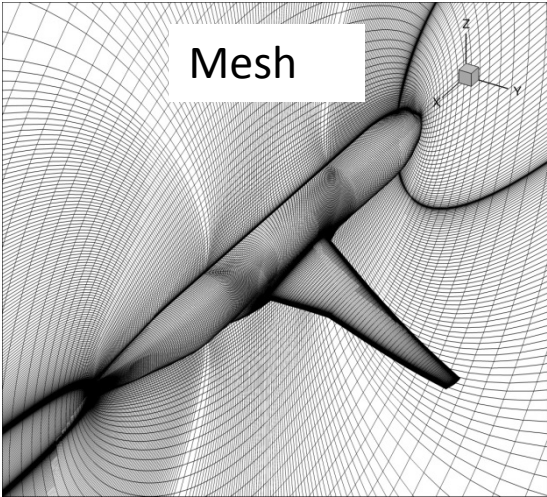
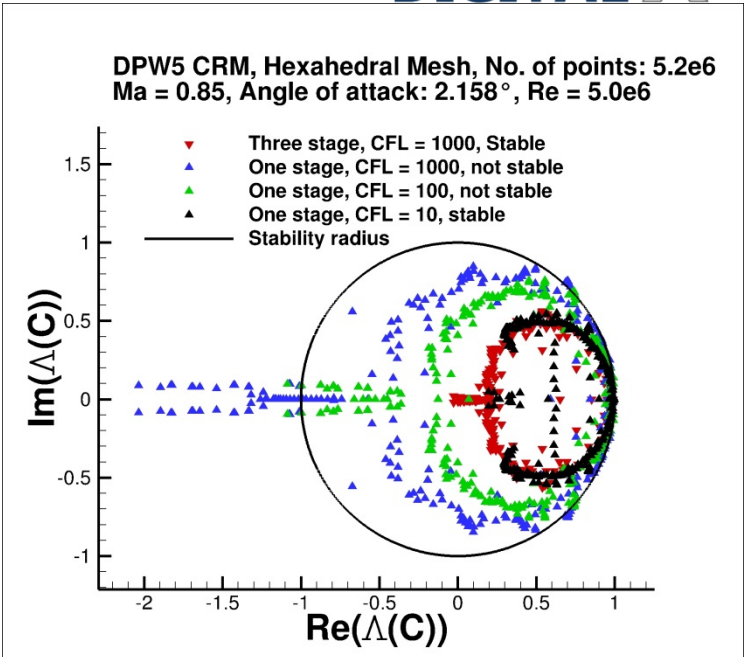
# Analysis of schemes: Impact of CFL number

## Analysis for mesh with 5.2e6 points:

### Investigation of number of stages for

- symmetric Line Gauss-Seidel
- different CFL numbers

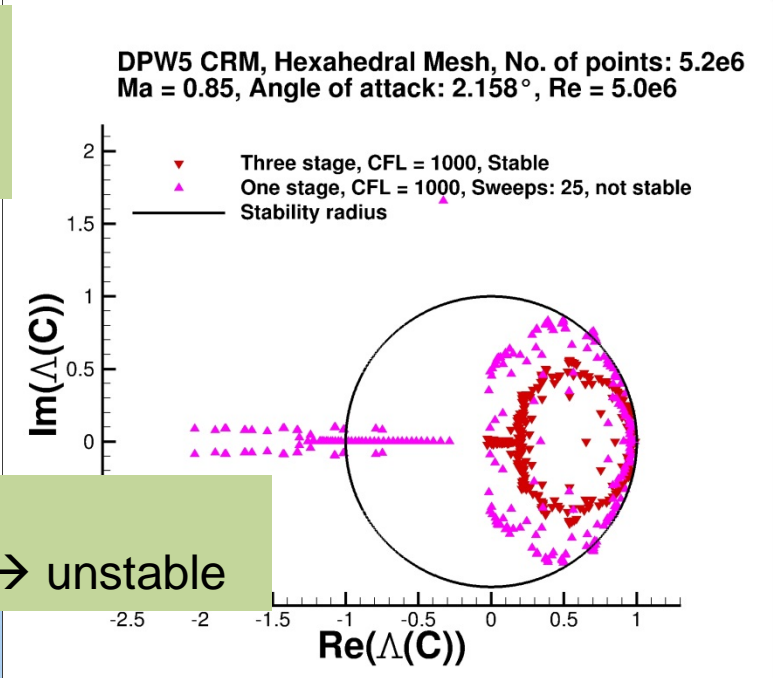
One stage: CFL = 1000 → unstable  
One stage: CFL = 100 → unstable  
One stage: CFL = 10 → stable  
Three stage: CFL = 1000 → stable



Significant reduction  
of CFL necessary for  
one stage schemes

Additional effort  
does not pay of

Sweeps: 25  
One stage: CFL = 1000 → unstable



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  - Spalart-Allmaras (neg.)
  - Wilcox ( $k\omega$ )



1992 model:

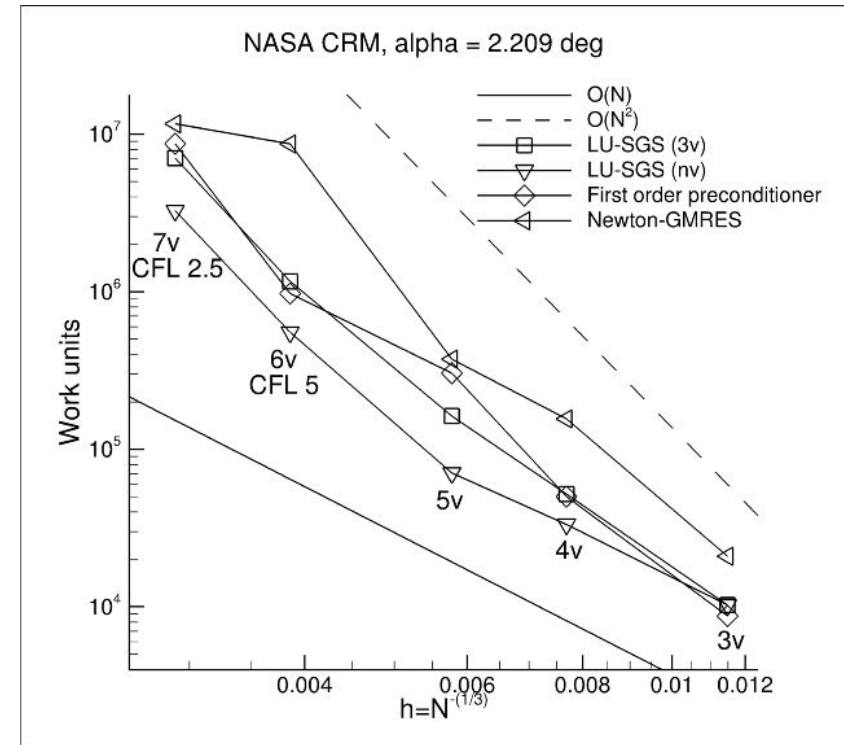
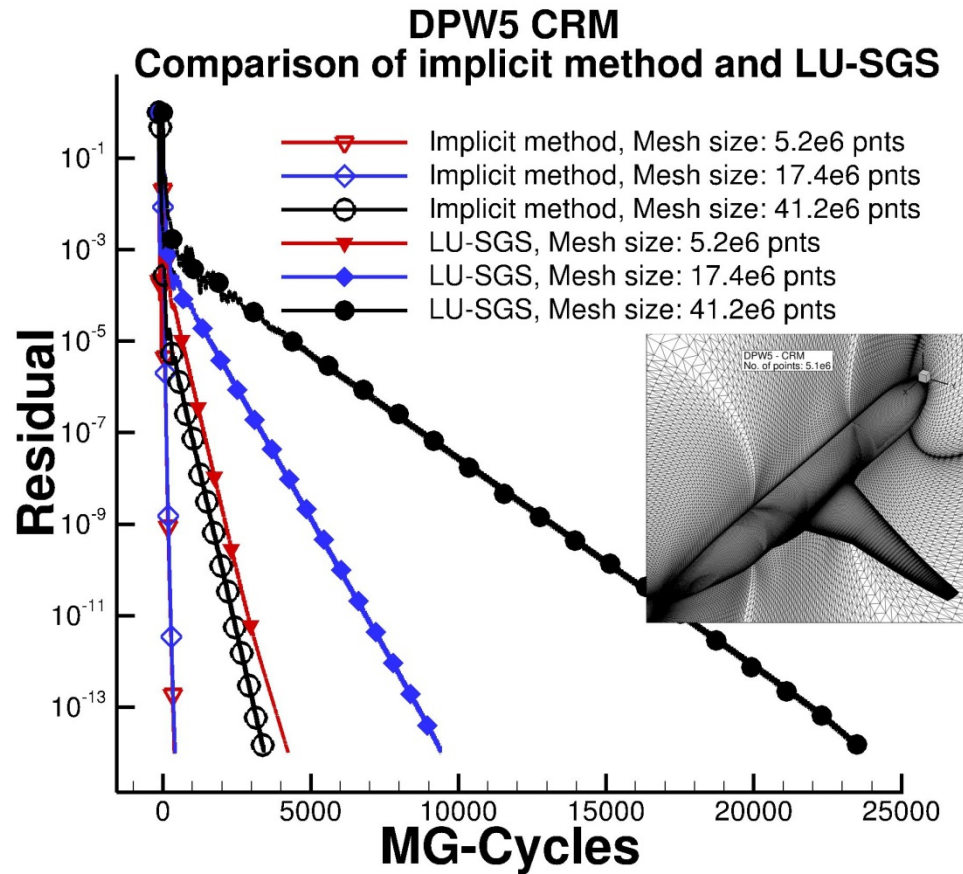
- transported variable can become negative such that iteration diverges
- choice of farfield values not clarified
- ....

2012 modification of original model:

- Allows for small negative values of transported variable
- clarification of choice of farfield values
- description and recommendations of implementation of several terms and details

→ 2012 version has been successfully implemented into the DLR TAU-Code





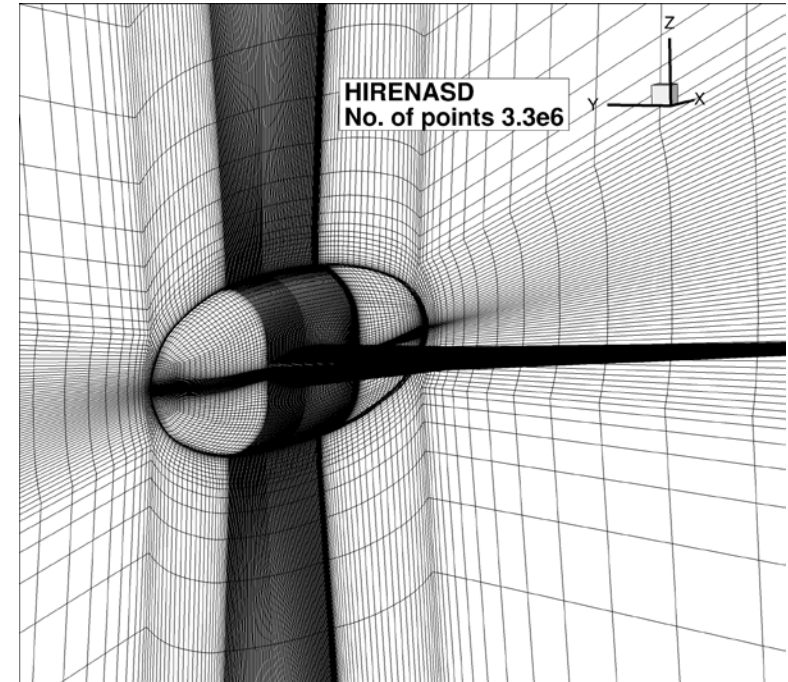
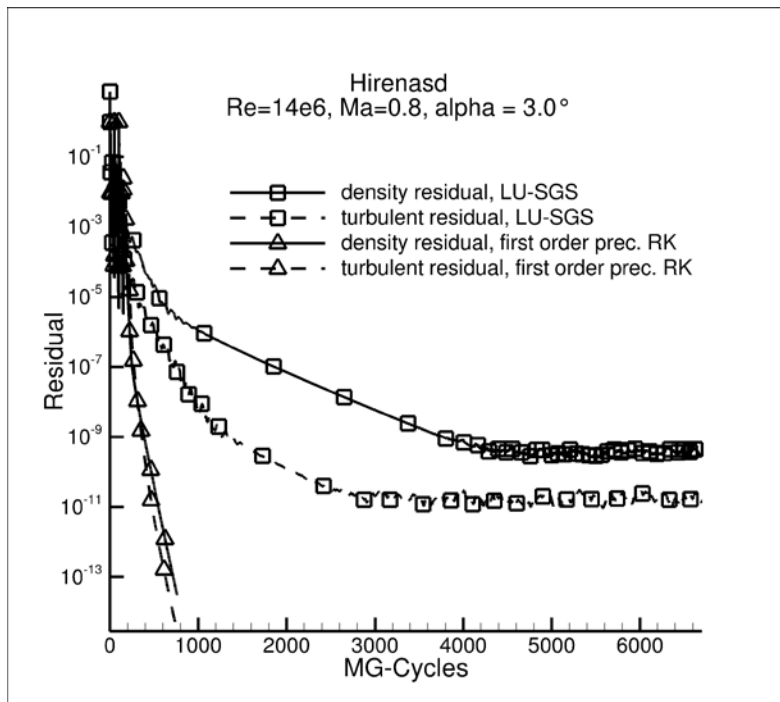
**Implicit methods have comparable complexity to standard LU-SGS method by improved robustness**





# Numerical example: HIRENASD

- High Reynolds number Aero-Structural Dynamics wind tunnel configuration
- $Ma = 0.8$
- $\alpha = 3.0^\circ$
- $Re = 14e6$
- Pure hexahedral mesh:  $3.3e6$  points



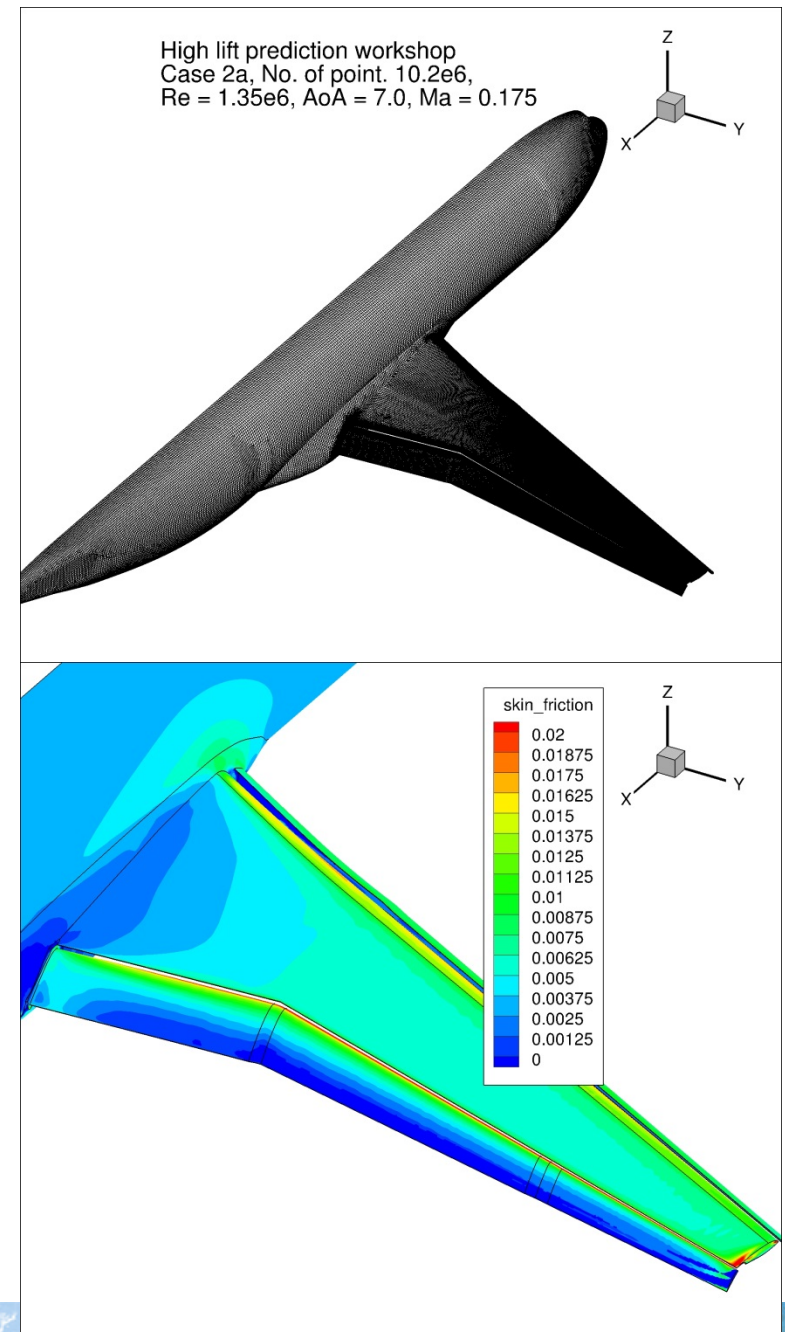
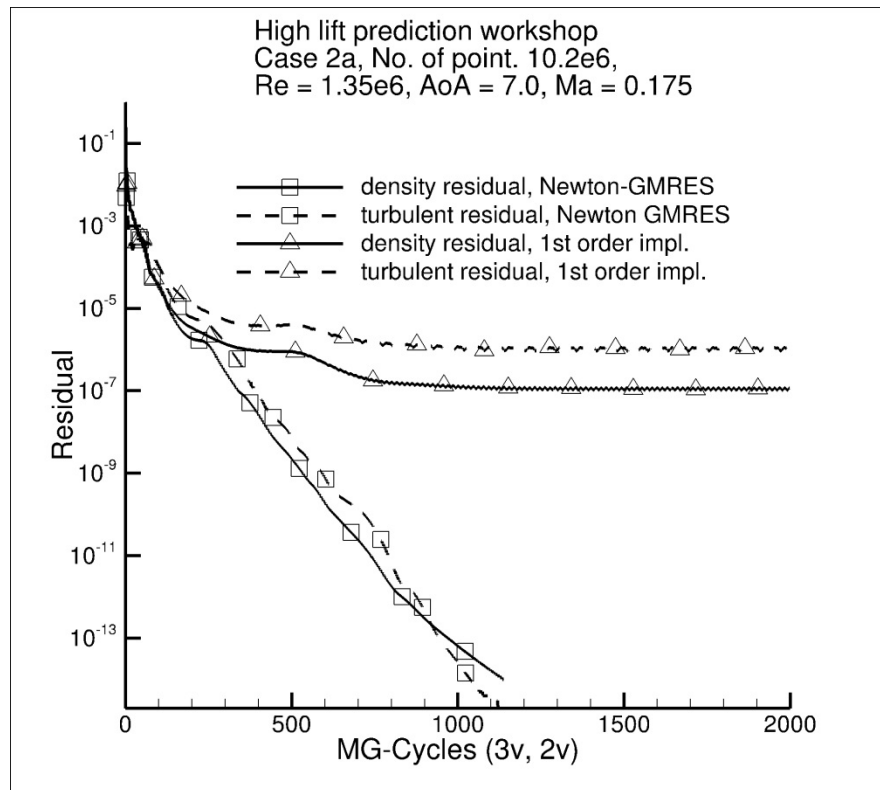
→ Implicit method converges,  
→ LU-SGS method stalls

# Configuration from second high-lift prediction workshop: Case 2a

## Necessity of Newton-kind algorithms

Turbulence model: SA-Neg

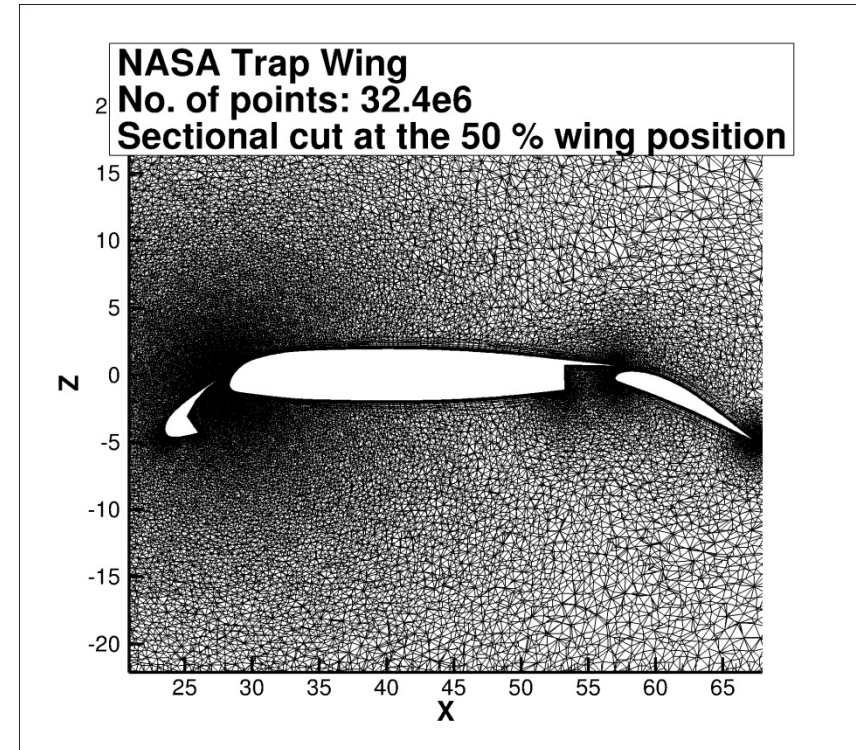
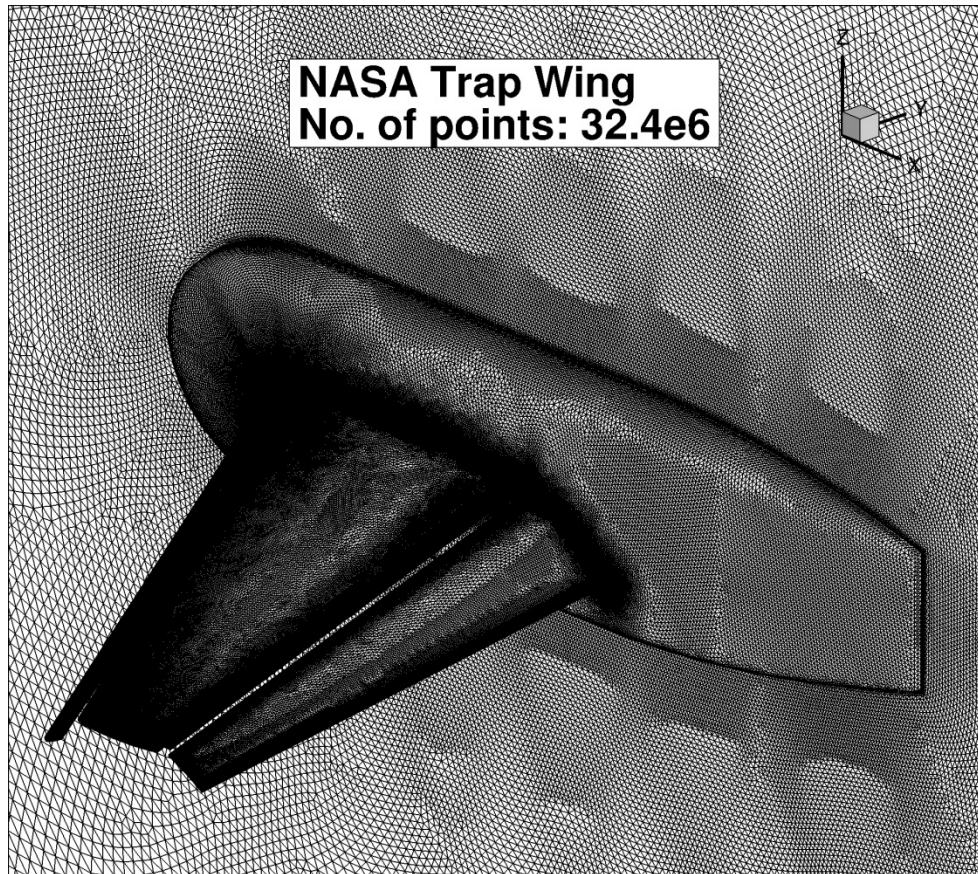
$Ma = 0.175$ ,  $Re = 1.35e6$ ,  $AoA = 7.0^\circ$





# NASA Trap Wing, $Ma = 0.2$ , $Re = 4.3e6$

Unstructured mesh results for  
 $AOA = 13^\circ, 28^\circ, 32^\circ, 34^\circ, 37^\circ$



- Coarse Mesh: 3.7e6 NDOF
- Medium Mesh: 11.0e6 NDOF
- Fine Mesh: 32.4e6 NDOF

VGRID Meshes used at High Lift Prediction Workshop 1

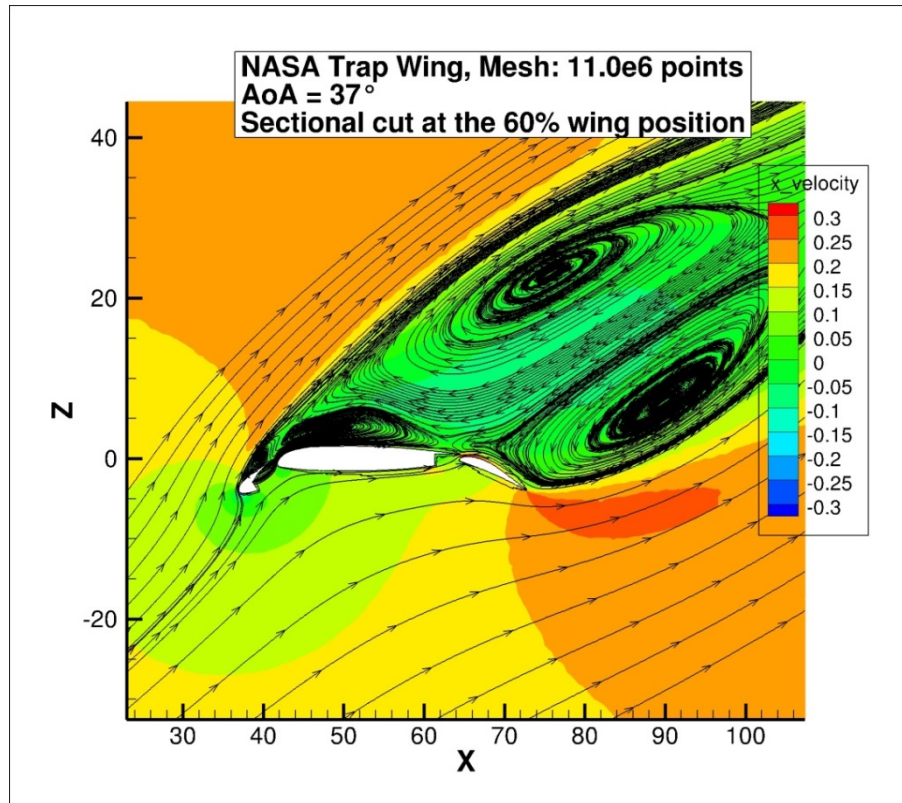




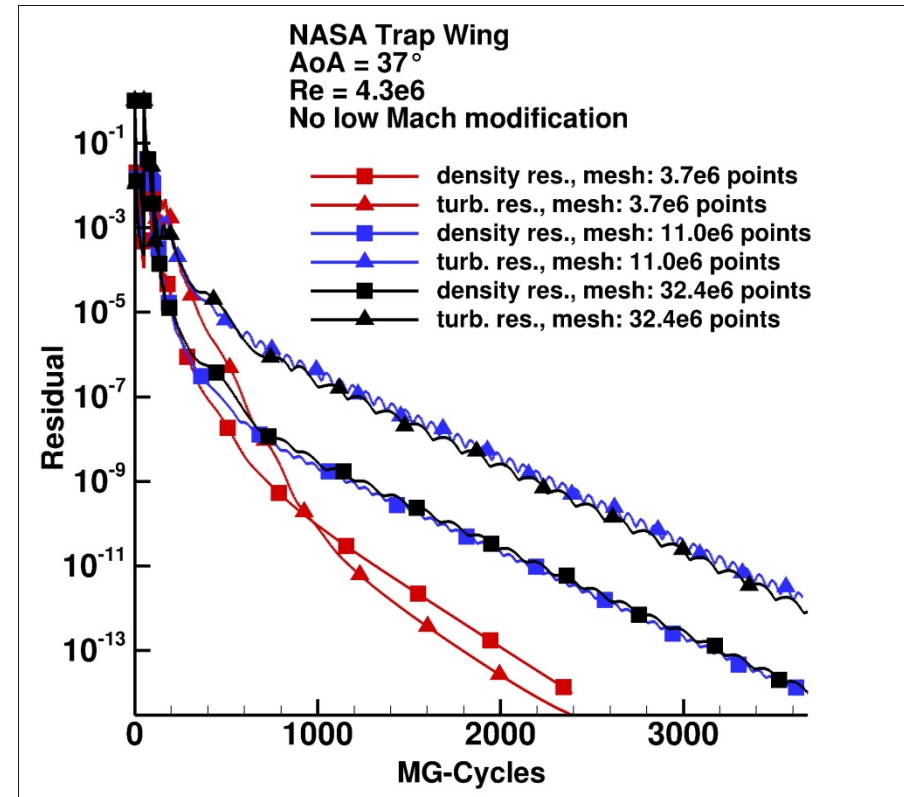
# NASA Trap Wing, $Ma = 0.2$ , $Re = 4.3e6$

Unstructured mesh results for  $AOA = 37^\circ$

- Residual has been reduced to machine accuracy using Newton kind methods  
→ Steady state could not be found with simplified algorithms



Flow field at the 60% wing section



Convergence history for  $AoA = 37^\circ$



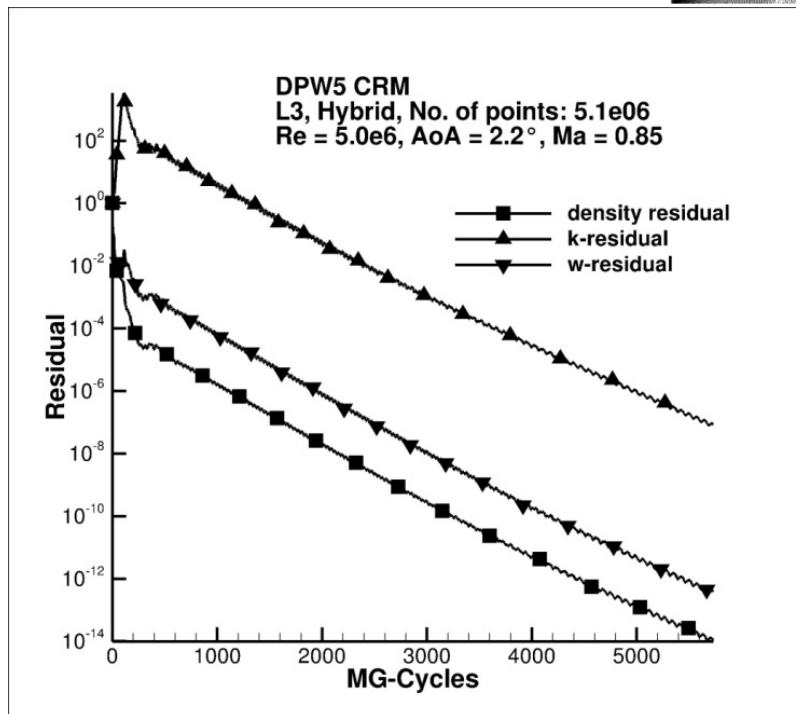
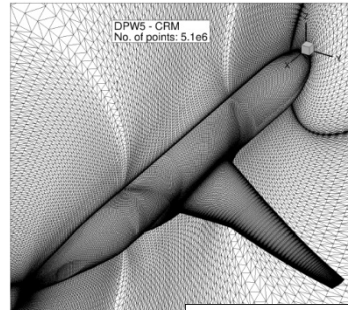
# Outline

- Introduction and Motivation – The RANS equations
- Solution algorithms – Multigrid smoothers
- Best practice considerations
- Numerical examples
  - Spalart-Allmaras (neg.)
  - Wilcox ( $k\omega$ )

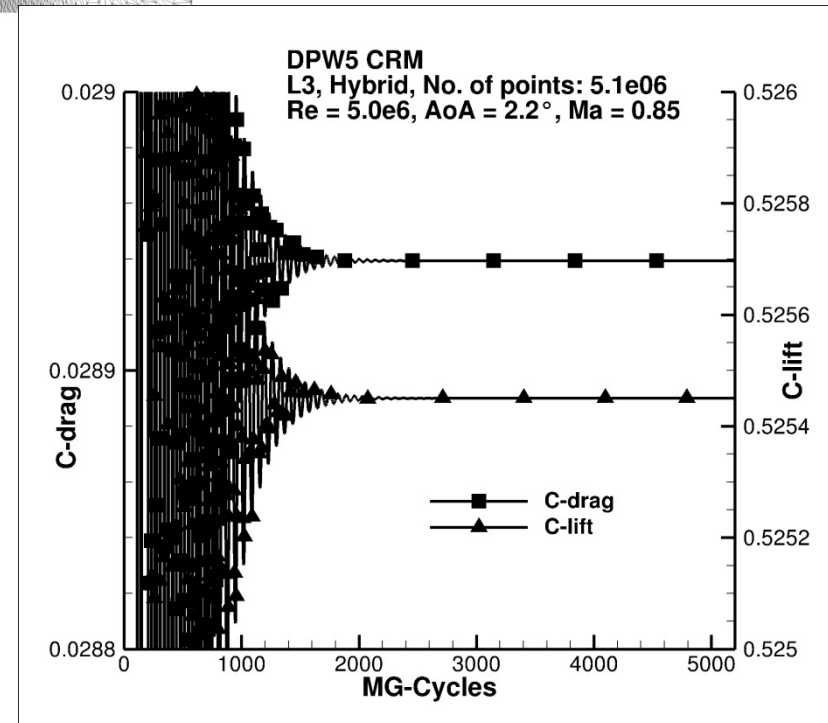


# Numerical examples: DPW5 + $k\omega$ -model

- Wing-body configuration
- $Ma = 0.85$
- $\alpha = 2.209^\circ$
- $Re = 5e6$
- No. of points:  $5.1e6$



Convergence history of residuals

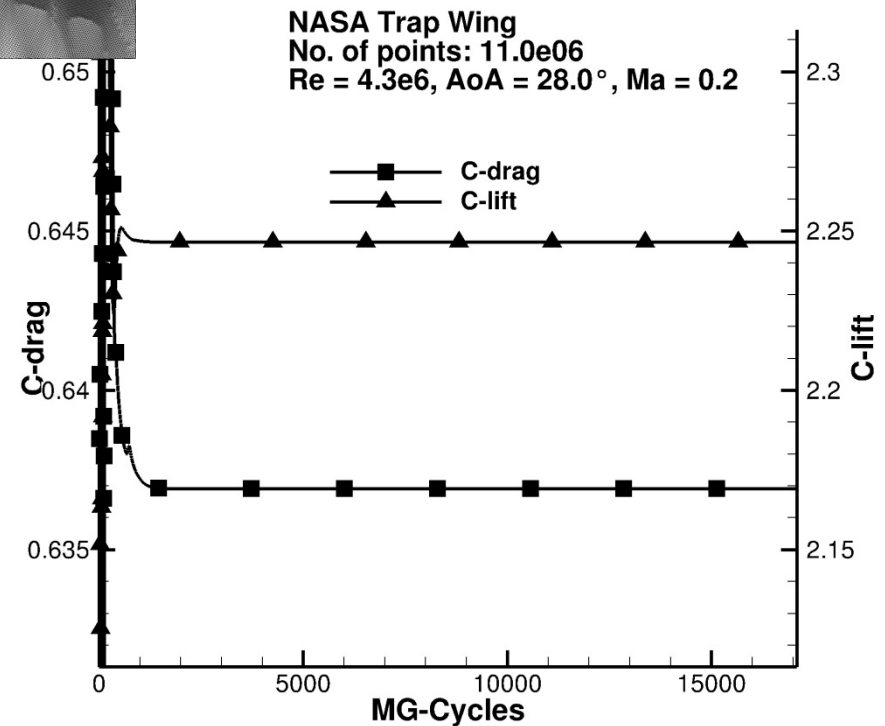
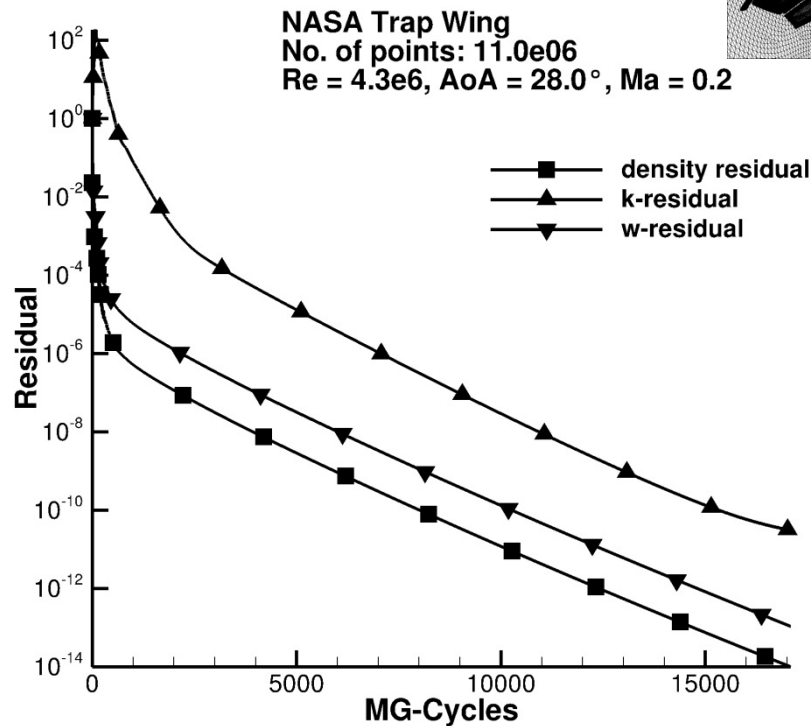
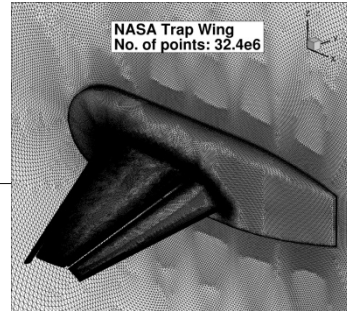


Convergence history of lift and drag



# NASA Trap Wing + $k\omega$ -model

$Ma = 0.2$ ,  $Re = 4.3e6$ ,  $AoA = 28.0^\circ$



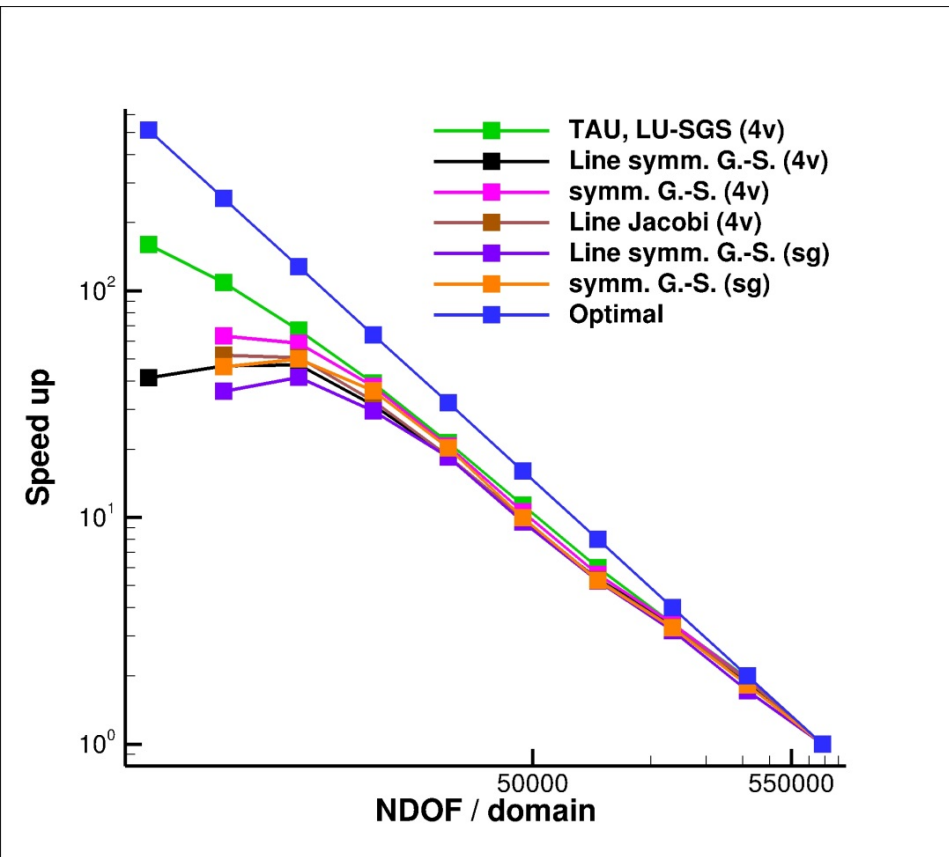
Convergence history of residuals

Convergence history of lift and drag

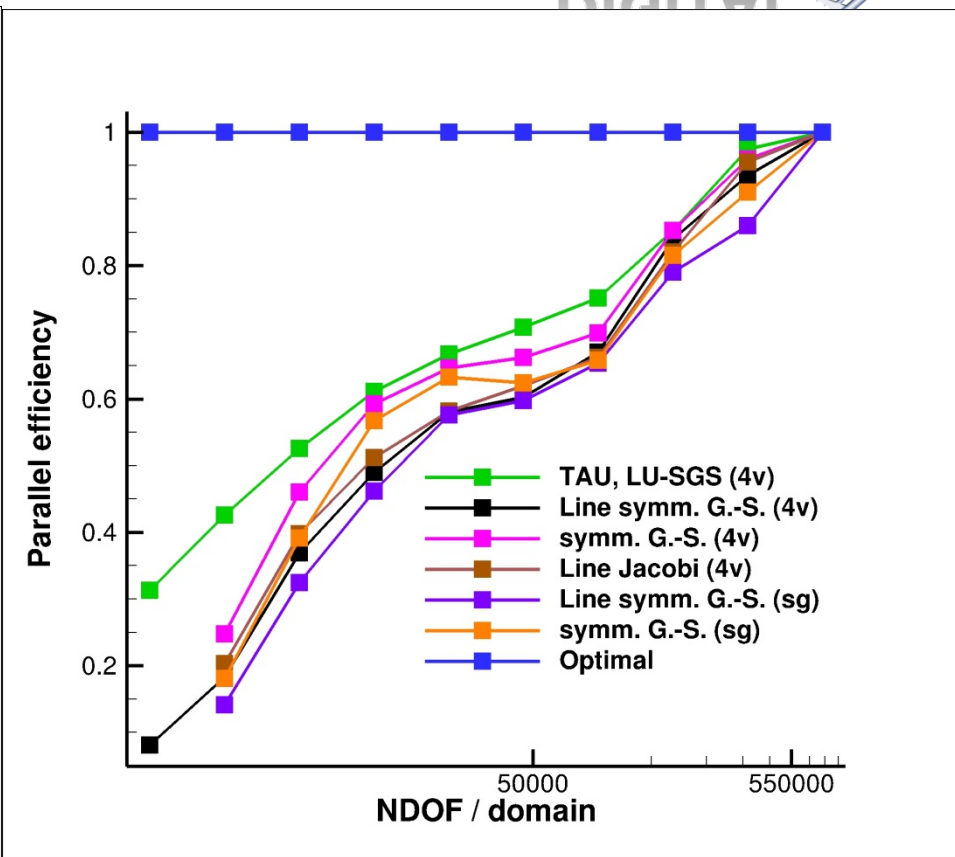




# Speed up and parallel efficiency: Strong scaling



Actual speed up



Actual parallel efficiency,  
System effectiveness

→ Severe issue with respect to exploitation of modern hardware clusters





# Summary

## Implicit methods offer the potential to

- improve significantly the observed convergence rates
- find fully (machine accurate) converged solutions of complex flows
- significantly increase robustness (e.g. they work for a broad range of CFL numbers)
- implement the hierarchy of smoothers in one framework
- outsource and decouple the main work into a suited linear algebra package

## Implicit methods require

- significantly more time per iteration than explicit methods
- a fully differentiated code which needs to be kept up to date
- significant more fast memory
- are not straightforward to ensure good parallel scalability
- to outsource and decouple the main work into a suited **linear algebra package**
- a new framework → **Flucs code**



Thank you!  
Questions?

